

-N-<sup>13</sup>C<sup>α</sup>-<sup>13</sup>C'-N-C<sup>α</sup>-C'-N-<sup>13</sup>C<sup>α</sup>-<sup>13</sup>C'-<sup>15</sup>N-C<sup>α</sup>-C'-N-C<sup>α</sup>-C'-  
Arg45 - Tyr46 - Arg47 - Asp48 - Val49

**Fig. 1**

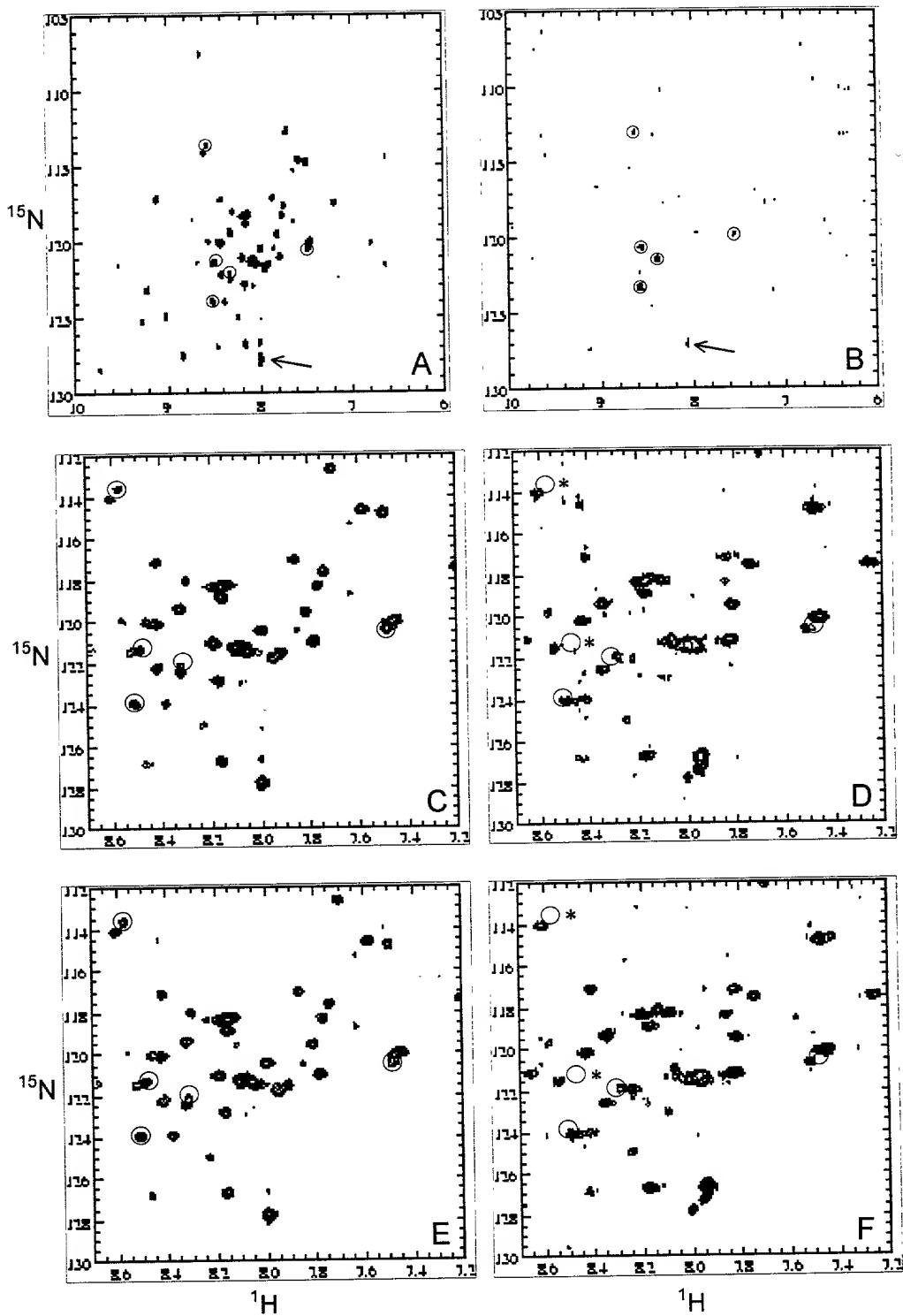


Fig. 2



**Fig. 3**

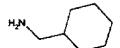
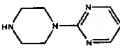
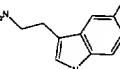
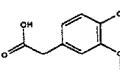
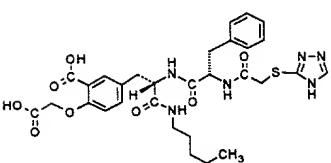
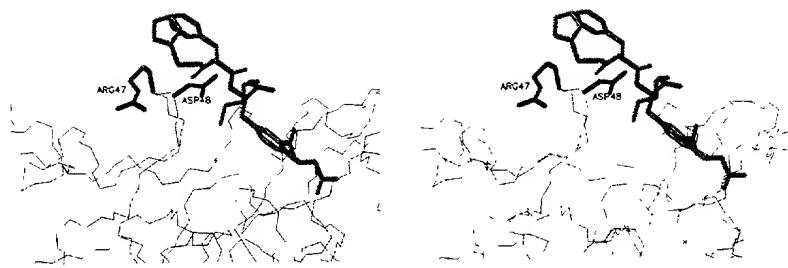
Compound	Structure	% inhibition	NMR binder
control (no compound added)		0	-
N35		14	no
N136		28	no
N200		20	yes
N212		0	no
PNU179983		100	yes

Fig. 4



**Fig. 5**

## Application I: screen for binding to single binding site

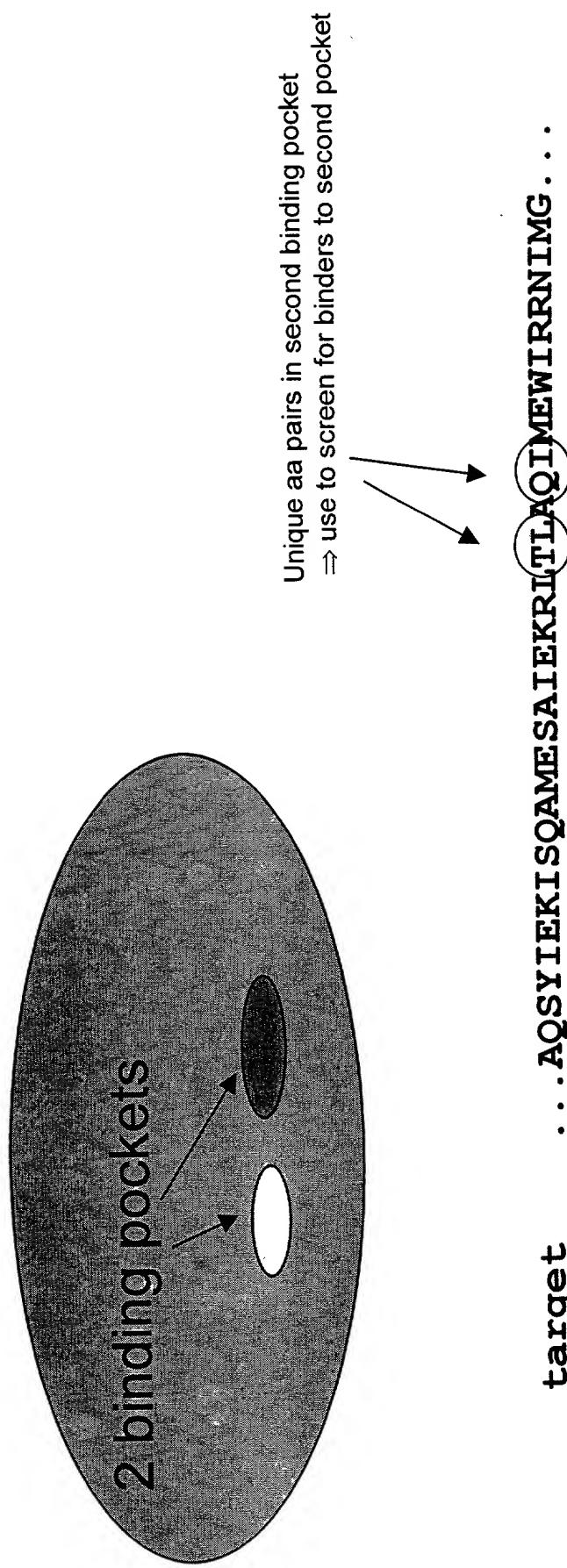


Fig. 6

## Application II: screen for specificity in vicinity of active site

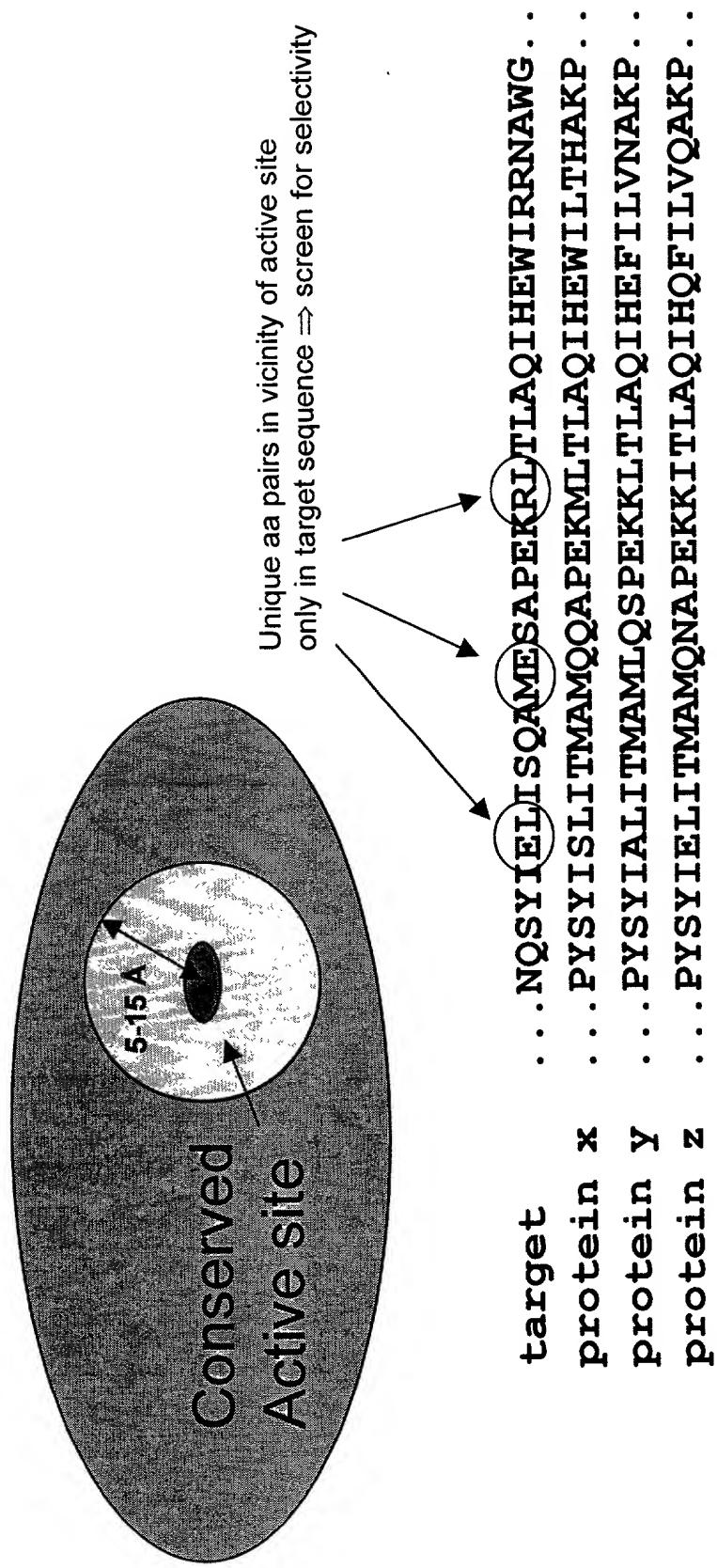


Fig. 7

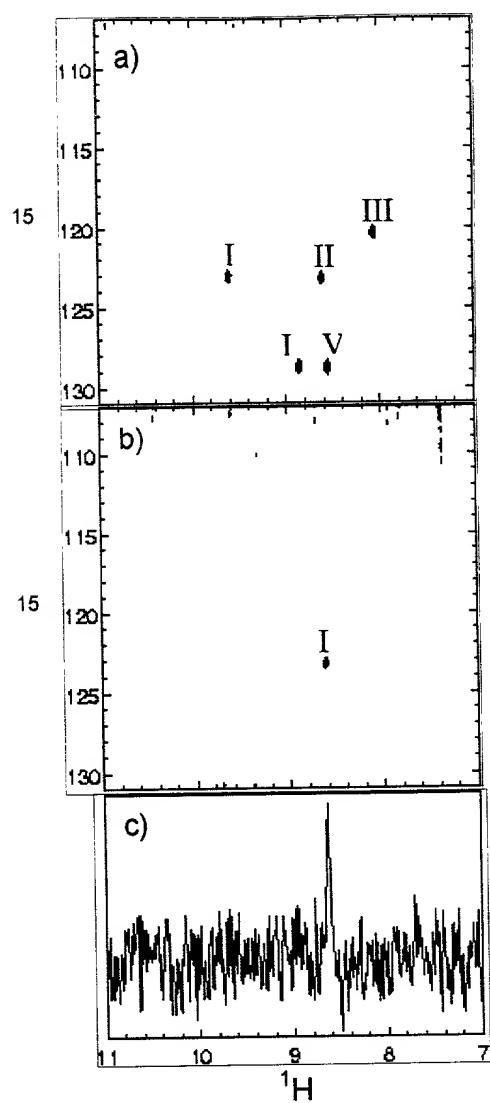
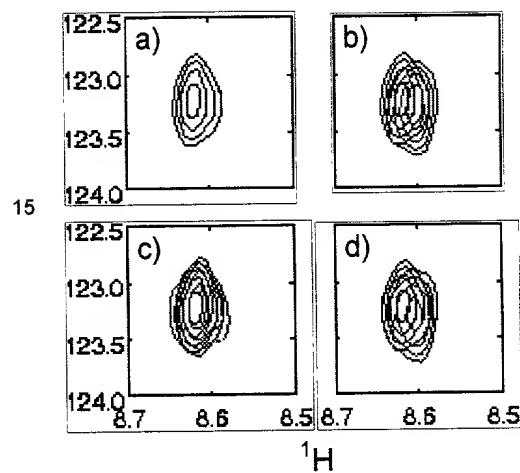


Fig. 8



**Fig. 9**